7411 Beach Drive E, Port Orchard Washington 98366

CASE NARRATIVE

January 29, 1999

Project:

Hamilton/Labree

Samples:

99048005

Project ID:

109699

Project Officer:

Pam Marti

By:

Karin Feddersen (

VOLATILE ORGANIC ANALYSIS

SUMMARY:

The data is usable as reported.

ANALYTICAL METHODS:

Volatile organic compounds were analyzed using the Manchester Laboratory modification of the EPA Method 8260 purge-trap procedure and capillary Gas Chromatography with Mass Spectrometer (GC/MS) analysis. Routine QA/QC procedures were performed.

BLANKS:

No analytes of interest were detected above the reporting limit in the blank.

SURROGATES:

Surrogate recoveries were within acceptable limits for all samples.

HOLDING TIMES:

The samples were analyzed within the recommended 14 day holding time.

MATRIX SPIKE AND MATRIX SPIKE DUPLICATE:

Insufficient sample was provided for matrix spikes. At least three extra vials are required to perform matrix spikes.



DATA QUALIFIER CODES:

U The analyte was not detected at or above the reported value. J The analyte was positively identified. The associated numerical value is an estimate. UJ The analyte was not detected at or above the reported estimated result. **REJ** The data are unusable for all purposes. NAF Not analyzed for. N There is evidence the analyte is present in the sample. NJ There is evidence that the analyte is present. The associated numerical result is an estimate. E This qualifier is used when the concentration of the associated value exceeds the known calibration range. The associated numerical result is an estimate.

bold - The analyte was present in the sample. (Visual Aid to locate detected compounds on report sheet.)

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Analysis Report for

Volatile Organic Analysis

Hamilton Labree Rds. - 04 Project Name:

LIMS Project ID: 1096-99

Sample: 99048005

Date Collected: 01/26/99

Method: SW8260 Matrix: Water

Field ID: HL-5

Project Officer: Pam Marti

Date Analyzed: 01/29/99

Units: ug/L

Analyte	Result	Qualifier	Analyte	Result	Qualifier
Dichlorodifluoromethane	20	U	Tetrachloroethene	7.1	
Chloromethane	5	U	Dibromochloromethane	2	U
Vinyl Chloride	1	U	1,2-Dibromoethane (EDB)	1	U
Bromomethane	1	U	Chlorobenzene	1	U
Chloroethane	5	U	1,1,1,2-Tetrachloroethane	2 5	U
Trichlorofluoromethane	1	U	Ethylbenzene	5	U
Ethyl Ether	1	U	m & p-Xylene	2	U
1,1,2 Trichlorotrifluoroethane	1	U	o-Xylene	1	U
1, 1-Dichloroethene	5	U	Styrene	2	U
Acetone	2	U	Bromoform	2	U
Methyl Iodide	1	U	Isopropylbenzene (Cumene)	1	U
Carbon Disulfide	2	U	1,1,2,2-Tetrachloroethane	1	U
Methylene Chloride	1	U	Trans-1,4-Dichloro-2-butene	5 2	U
2-Methyoxy-2-Methylpropane	2	U	1,2,3-Trichloropropane	2	U
Trans-1,2-Dichloroethene	1	U	Bromobenzene	1	U
1,1-Dichloroethane	1	U	n-Propylbenzene	1	U
2-Butanone	5	U	2-Chlorotoluene	2 2 2 2 2	U
Cis-1,2-Dichloroethene	1	U	1,3,5-Trimethylbenzene	2	U
2,2-Dichloropropane	2	U	4-Chlorotoluene	2	U
Bromochloromethane	1	U	Tert-Butylbenzene	2	U
Chloroform	1	U	1,2,4-Trimethylbenzene	2	U
Tetrahydrofuran	2	U	Pentachloroethane	1	U
1,1,1-Trichloroethane	1	U	Sec-Butylbenzene	1	U
1,1-Dichloropropene	10	U	p-Isopropyltoluene	2	ប
Carbon Tetrachloride	1	U	1,3-Dichlorobenzene	2	U
1,2-Dichloroethane	1	U	1,4-Dichlorobenzene	2	U
Benzene	5	U	n-Butylbenzene	5	Ū
Trichloroethene	1	U	1.2-Dichlorobenzene	1	U
1,2-Dichloropropane	5	U	Hexachloroethane	5	U
Dibromomethane	5	U	1,2-Dibromo-3-Chloropropane	5 5 5 2 5	Ū
Bromodichloromethane	2	Ū	1,2,4-Trichlorobenzene	5	Ū
Cis-1,3-Dichloropropene	2.1	Ū	Hexachlorobutadiene	2	Ū
4-Methyl-2-Pentanone	2	Ū	Naphthalene	5	Ü
Toluene	1	Ŭ	1,2,3-Trichlorobenzene	5	Ŭ
Trans-1,3-Dichloropropene	ī.9	Ŭ		_	_
1,1,2-Trichloroethane	2	Ŭ			
1,3-Dichloropropane	$\bar{2}$	Ŭ			
2-Hexanone	10	Ŭ			

Authorized By:

Release Date: ___2/5/99

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Analysis Report for

Volatile Organic Analysis

Project Name: Hamilton Labree Rds. - 04 LIMS Project ID: 1096-99

Sample: 99048005

Date Collected: 01/26/99

Method: SW8260

Field ID: HL-5

Matrix: Water

Project Officer: Pam Marti

Date Analyzed: 01/29/99

Units:

ug/L

Surrogate Recoveries

1,2-Dichloroethane-D4	102	%
1,4-Difluorobenzene	112	%
Toluene-D8	110	%
p-Bromofluorobenzene	97	%
1,2-Dichlorobenzene-D4	104	%

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Analysis Report for

Volatile Organic Analysis

Project Name: Hamilton Labree Rds. - 04 LIMS Project ID: 1096-99

ODBW9029 Lab ID:

Method: SW8260

QC Type: Laboratory Method Blank Project Officer: Pam Marti

Matrix: Water

Date Analyzed: 01/29/99

Units: ug/L

Analyte	Result	Qualifier	Analyte	Result	Qualifier
Dichlorodifluoromethane	20	U	Tetrachloroethene	1	U
Chloromethane	5	U	Dibromochloromethane	2	U
Vinyl Chloride	1	U	1,2-Dibromoethane (EDB)	1	U
Bromomethane	1	U	Chlorobenzene	1	U
Chloroethane	5	U	1,1,1,2-Tetrachloroethane	2 5 2	U
Trichlorofluoromethane	1	U	Ethylbenzene	5	U
Ethyl Ether	1	U	m & p-Xylene		U
1,1,2 Trichlorotrifluoroethane	1	U	o-Xylene	1	U
1,1-Dichloroethene	5	U	Styrene	2 2 1	U
Acetone	2	U	Bromoform	2	${f u}$
Methyl Iodide	1	U	Isopropylbenzene (Cumene)	1	U
Carbon Disulfide	2	U	1,1,2,2-Tetrachloroethane	1 5 2	U
Methylene Chloride	1	U	Trans-1,4-Dichloro-2-butene	5	U
2-Methyoxy-2-Methylpropane	2	U	1,2,3-Trichloropropane	2	U
Trans-1,2-Dichloroethene	1	U	Bromobenzene	1	U
1,1-Dichloroethane	1	U	n-Propylbenzene	1	U
2-Butanone	5	U	2-Chlorotoluene	2	U
Cis-1,2-Dichloroethene	1.	U	1,3,5-Trimethylbenzene	2	U
2,2-Dichloropropane	2	U	4-Chlorotoluene	2 2 2 2 2	U
Bromochloromethane	1	U	Tert-Butylbenzene	2	U
Chloroform	1	U	1,2,4-Trimethylbenzene	2	U
Tetrahydrofuran	2	U	Pentachloroethane	1	U
1,1,1-Trichloroethane	1	U	Sec-Butylbenzene	1	U
1,1-Dichloropropene	10	U	p-Isopropyltoluene	2	U
Carbon Tetrachloride	1	U	1,3-Dichlorobenzene	2 2 1	U
1,2-Dichloroethane	1	Ū	1,4-Dichlorobenzene	1	บ
Benzene	5	Ū	n-Butylbenzene	5	U
Trichloroethene	1	Ū	1,2-Dichlorobenzene	1	Ŭ
1,2-Dichloropropane	5	Ū	Hexachloroethane	5	Ū
Dibromomethane	5	Ŭ	1,2-Dibromo-3-Chloropropane	5	Ū
Bromodichloromethane	2	$\check{\mathbf{U}}$	1,2,4-Trichlorobenzene	5	Ŭ
Cis-1,3-Dichloropropene	2 2.1	Ŭ	Hexachlorobutadiene	5 5 5 2 5	Ŭ
4-Methyl-2-Pentanone	2	Ŭ	Naphthalene	5	Ŭ
Toluene	1	Ŭ	1,2,3-Trichlorobenzene	5	Ŭ
Trans-1,3-Dichloropropene	1.9	Ŭ	- 1-10 TIONIOLOUININ		~
1,1,2-Trichloroethane		ប័			
1,3-Dichloropropane	2 2	บี			
2-Hexanone	10	Ŭ			
2 Horanono	10	O			

Authorized By:

Release Date: 2/5/99

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Analysis Report for

Volatile Organic Analysis

Project Name:

Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Method: SW8260

Lab ID: ODBW9029 QC Type: Laboratory Method Blank Project Officer: Pam Marti

Date Analyzed: 01/29/99

Matrix: Water Units: ug/L

Surrogate Recoveries

1,2-Dichloroethane-D4	99	%
1,4-Difluorobenzene	110	%
Toluene-D8	106	%
p-Bromofluorobenzene	90	%
1,2-Dichlorobenzene-D4	105	%

Authorized By:

Release Date: 2/5/99

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MANCHESTER ENVIRONMENTAL LABORATORY

7411 Beach Drive E, Port Orchard Washington 98366

February 5, 1999

Subject:

Hamilton Labree

Samples:

99048005

Project ID:

109699

Project Officer:

Pam Marti

By:

Karin Feddersen

SEMIVOLATILE ORGANICS

ANALYTICAL METHODS:

The samples were extracted following the EPA CLP and SW-846 8270 procedure. The extracts were cleaned up with Gel Permeation Chromatography (GPC). Analysis was by capillary gas chromatography with mass spectrometry (GC/MS). Routine QA/QC procedures were performed with the analyses.

HOLDING TIMES:

The samples were stored at 4 degrees C until extraction. They were extracted and analyzed within the recommended holding times.

BLANKS:

Low levels of some analytes were detected in the laboratory blanks. An analyte is considered native to the sample when the on-column concentration is at least five times greater than in the associated method blanks. A phthalate is considered native to the sample when the concentration is at least ten times greater than in the associated method blanks.

SURROGATES:

The standard Manchester Laboratory Base/Neutral/Acid (BNA) surrogates were added to the sample prior to extraction. All surrogate recoveries were within acceptable limits.

MATRIX SPIKE AND MATRIX SPIKE DUPLICATE:

Insufficient sample was provided for spikes.

COMMENTS:

The data is acceptable for use as reported.

DATA QUALIFIER CODES:

The analyte was not detected at or above the reported value.

J - The analyte was positively identified. The associated numerical value is an estimate.

UJ - The analyte was not detected at or above the reported estimated result.

REJ - The data are unusable for all purposes.

NAF - Not analyzed for.

N - There is evidence the analyte is present in the sample.

NJ - There is evidence that the analyte is present. The associated numerical result is an estimate.

E - This qualifier is used when the concentration of the associated value exceeds the known calibration range. The associated numerical result is an estimate.

bold - The analyte was present in the sample. (Visual Aid to locate detected compounds on report sheet.)

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Analysis Report for

Base/Neutral/Acids

Project Name: Hamilton Labree Rds. - 04 LIMS Project ID: 1096-99

Sample: 99048005 Date Collected: 01/26/99

Method: SW8270

Field ID: HL-5

Date Prepared: 01/27/99

Matrix: Water

Project Officer: Pam Marti

Date Analyzed: 01/27/99 Units: ug/L

Analyte	Result	Qualifier	Analyte	Result	Qualifier
N-Nitrosodimethylamine	.25	U	Acenaphthene	.12	ប
Pyridine	.12	U	2,4-Dinitrophenol	2.5	ប
Aniline	.12	U	4-Nitrophenol	.12	U
Phenol	.12	U	Dibenzofuran	.12	U
Bis(2-Chloroethyl)Ether	.12	U	2,4-Dinitrotoluene	.12	U
2-Chlorophenol	.12	U	Diethylphthalate	.12	U
1,3-Dichlorobenzene	.12	U	Fluorene	.12	U
1,4-Dichlorobenzene	.12	U	4-Chlorophenyl-Phenylether	.12	U
1,2-Dichlorobenzene	.12	U	4-Nitroaniline	.12	U
Benzyl Alcohol	.12	Ü	4,6-Dinitro-2-Methylphenol	1.2	U
2-Methylphenol	.12	U	N-Nitrosodiphenylamine	.12	U
2,2'-Oxybis[1-chloropropane]	.12	Ū	1,2-Diphenylhydrazine	.12	U
N-Nitroso-Di-N-Propylamine	.12	Ū	4-Bromophenyl-Phenylether	.12	U
4-Methylphenol	.12	U	Hexachlorobenzene	.12	U
Hexachloroethane	.12	Ū	Pentachlorophenol	.25	U
Nitrobenzene	.12	Ū	Phenanthrene	.12	Ū
Isophorone	.12	Ū	Anthracene	,12	Ū
2-Nitrophenol	.12	Ū	Caffeine	.12	Ū
2,4-Dimethylphenol	.12	Ū	Carbazole	.12	Ū
Bis(2-Chloroethoxy)Methane	.12	Ū	Di-N-Butylphthalate	.12	Ŭ
Benzoic Acid		REJ	Fluoranthene	.12	Ŭ
2,4-Dichlorophenol	.12	Ū	Benzidine	.25	Ŭ
1,2,4-Trichlorobenzene	.12	Ŭ	Pyrene	.12	Ŭ
Naphthalene	.12	Ŭ	Retene	.12	Ü
4-Chloroaniline	.12	Ŭ	Butylbenzylphthalate	.12	Ŭ
Hexachlorobutadiene	.12	Ŭ	Benzo(a)anthracene	.12	Ŭ
4-Chloro-3-Methylphenol	.12	Ŭ	3,3'-Dichlorobenzidine	.25	Ŭ
2-Methylnaphthalene	.12	Ŭ	Chrysene	.12	Ŭ
1-Methylnaphthalene	.12	Ŭ	Bis(2-Ethylhexyl) Phthalate	.12	Ŭ
Hexachlorocyclopentadiene	.12	Ŭ	Di-N-Octyl Phthalate	.12	Ŭ
2 4 6-Trichlorophenol	12	Ŭ	Benzo(b)fluoranthene	.12	ប័
2,4,5-Trichlorophenol	.12	Ŭ	Benzo(k)fluoranthene	.12	Ŭ
2-Chloronaphthalene	.12	บั	Benzo(a)pyrene	.12	Ŭ
2-Nitroaniline	.12	ŭ	3B-Coprostanol	.25	Ŭ
Dimethylphthalate	.12	Ŭ	Indeno(1,2,3-cd)pyrene	.12	ŭ
2,6-Dinitrotoluene	.12	Ŭ	Dibenzo(a,h)anthracene	.12	Ŭ
Acenaphthylene	.12	Ŭ	Benzo(ghi)perylene	.12	Ŭ
3-Nitroaniline	.12	Ŭ	Dougo (Bur) por Jiono	.12	J
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Authorized By:

Release Date: 2/5/99

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Analysis Report for

Base/Neutral/Acids

Project Name: Hamilton Labree Rds. - 04 LIMS Project ID: 1096-99

Sample: 99048005

Date Collected: 01/26/99

Method: SW8270

Field ID: HL-5

Date Prepared: 01/27/99 Date Analyzed: 01/27/99

Matrix: Water

Project Officer: Pam Marti

Units:

ug/L

Surrogate Recoveries

2-Fluorophenol	39	%
D5-Phenol	22	%
D4-2-Chlorophenol	80	%
1,2-Dichlorobenzene-D4	66	%
D5-Nitrobenzene	88	%
2-Fluorobiphenyl	73	%
Pyrene-D10	119	%
Terphenyl-D14	114	%

Authorized By:

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Analysis Report for

Base/Neutral/Acids

Project Name: Hamilton Labree Rds. - 04 LIMS Project ID: 1096-99

Sample: 99048005

Date Collected: 01/26/99

Method: SW8270

Field ID: HL-5

Project Officer: Pam Marti

Date Prepared: 01/27/99 Date Analyzed: 01/27/99

Matrix: Water Units: ug/L

Tentatively Identified Compounds

CAS Number Analyte Description

Result Qualifier

127184

Tetrachloroethene

1.5

NJ

Authorized By: Release Date: 2/5/98

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Analysis Report for

Base/Neutral/Acids

Project Name: Hamilton Labree Rds. - 04 LIMS Project ID: 1/96-99

OBW9027B1 Lab ID:

QC Type: Laboratory Method Blank

Date Prepared: 01/27/99 Project Officer: Pam Marti

Method: SW8270 Matrix: Water

Date Analyzed: 01/27/99 Units: ug/L

Analyte	Result	Qualifier	Analyte	Result	Qualifier
N-Nitrosodimethylamine	.8	U	Acenaphthene	.4	U
Pyridine	.4	U	2,4-Dinitrophenol	8	U
Aniline	.4	U	4-Nitrophenol	.4	U
Phenol	.4	U	Dibenzofuran	.4	U
Bis(2-Chloroethyl)Ether	.4	U	2,4-Dinitrotoluene	.4	U
2-Chlorophenol	.4	U	Diethylphthalate	.13	J
1,3-Dichlorobenzene	.4	U	Fluorene	.4	U
1,4-Dichlorobenzene	.4	U	4-Chlorophenyl-Phenylether	.4	U
1,2-Dichlorobenzene	.4	U	4-Nitroaniline	,4	U
Benzyl Alcohol	.4	U	4,6-Dinitro-2-Methylphenol	4	U
2-Methylphenol	.4	U	N-Nitrosodiphenylamine	.4	U
2,2'-Oxybis[1-chloropropane]	.4	U	1,2-Diphenylhydrazine	.4	U
N-Nitroso-Di-N-Propylamine	.4	U	4-Bromophenyl-Phenylether	.4	U
4-Methylphenol	.4 .4	U	Hexachlorobenzene	.4	U
Hexachloroethane	.4	U	Pentachlorophenol	.8	U
Nitrobenzene	.4	U	Phenanthrene	.0043	J
Isophorone	.4	U	Anthracene	.4	U
2-Nitrophenol	.4	U	Caffeine	.4	U
2,4-Dimethylphenol	.4	U	Carbazole	.4	U
Bis(2-Chloroethoxy)Methane	.4	U	Di-N-Butylphthalate	.044	J
Benzoic Acid		REJ	Fluoranthene	.4	Ū
2,4-Dichlorophenol	.4	U	Benzidine	.8	U
1,2,4-Trichlorobenzene	.4	U	Pyrene	.4	U
Naphthalene	.4	U	Retene	.4	U
4-Chloroaniline	.4	Ū	Butylbenzylphthalate	.022	${f J}$
Hexachlorobutadiene	.4	Ū	Benzo(a)anthracene	.4	Ŭ
4-Chloro-3-Methylphenol	.4	Ū	3,3'-Dichlorobenzidine	.8	
2-Methylnaphthalene	.4	Ū	Chrysene	.4	Ū
1-Methylnaphthalene	.4	Ŭ	Bis(2-Ethylhexyl) Phthalate	.044	U J
Hexachlorocyclopentadiene	.4	Ŭ	Di-N-Octyl Phthalate	.4	ŭ
2,4,6-Trichlorophenol	.4	Ŭ	Benzo(b)fluoranthene	.4	Ŭ
2,4,5-Trichlorophenol	.4	Ŭ	Benzo(k)fluoranthene	.4	Ŭ
2-Chloronaphthalene	.4	Ŭ	Benzo(a)pyrene	A	Ŭ
2-Nitroaniline	.4	Ŭ	3B-Coprostanol	.8	Ŭ
Dimethylphthalate	.4	Ŭ	Indeno(1,2,3-cd)pyrene	.4	Ŭ
2,6-Dinitrotoluene	.4	Ū	Dibenzo(a,h)anthracene	.4	Ū
Acenaphthylene	.4	Ŭ	Benzo(ghi)perylene	.4	Ŭ
3-Nitroaniline	.4	Ŭ	(D) k) voice	- •	
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Analysis Report for

Base/Neutral/Acids

Project Name: Hamilton Labree Rds. - 04 LIMS Project ID: 1096-99

Lab ID: OBW9027B1

Date Prepared: 01/27/99

Method: SW8270

QC Type: Laboratory Method Blank Project Officer: Pam Marti

Date Analyzed: 01/27/99

Matrix: Water Units: ug/L

Surrogate Recoveries

2-Fluorophenol	66	%	
D5-Phenol	51	%	
D4-2-Chlorophenol	82	%	
1,2-Dichlorobenzene-D4	63	%	
D5-Nitrobenzene	83	%	
2-Fluorobiphenyl	76	%	
Pyrene-D10	110	%	
Terphenyl-D14	111	%	

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Analysis Report for

Base/Neutral/Acids

Hamilton Labree Rds. - 04 Project Name:

LIMS Project ID: 1096-99

OBW9027B2 Lab ID:

QC Type: Laboratory Method Blank Project Officer: Pam Marti Date Prepared: 01/27/99 Date Analyzed: 01/27/99

Method: SW8270 Matrix: Water

Units: ug/L

Analyte	Result	Qualifier	Analyte	Result	Qualifier
N-Nitrosodimethylamine	.8	U	Acenaphthene	.4	U
Pyridine	.4	U	2,4-Dînitrophenol	8	U
Aniline	.4	U	4-Nitrophenol	.4	U
Phenol	.4	U	Dibenzofuran	.4	U
Bis(2-Chloroethyl)Ether	.4	U	2,4-Dinitrotoluene	.4	U
2-Chlorophenol	.4	U	Diethylphthalate	.15	J
1,3-Dichlorobenzene	.4	U	Fluorene	.4	U
1,4-Dichlorobenzene	.4	U	4-Chlorophenyl-Phenylether	.4	U
1,2-Dichlorobenzene	.4	U	4-Nitroaniline	.4	U
Benzyl Alcohol	.4	U	4,6-Dinitro-2-Methylphenol	4	U
2-Methylphenol	.4	U ·	N-Nitrosodiphenylamine	.4 .4	U
2,2'-Oxybis[1-chloropropane]	.4	U	1,2-Diphenylhydrazine	.4	U
N-Nitroso-Di-N-Propylamine	4	U	4-Bromophenyl-Phenylether	.4	U
4-Methylphenol	.4 .4 .4	U	Hexachlorobenzene	.4	U
Hexachloroethane	.4	U	Pentachlorophenol	.8	U
Nitrobenzene	.4	U	Phenanthrene	.4	U
Isophorone	.4	U	Anthracene	.4 .4	U
2-Nitrophenol	.4	U	Caffeine	.4	U
2,4-Dimethylphenol	.4	U	Carbazole	.4	U
Bis(2-Chloroethoxy)Methane	.4	U	Di-N-Butylphthalate	.056	J
Benzoic Acid	16	U	Fluoranthene	,4	U
2,4-Dichlorophenol	,4	U	Benzidine	.8	U
1,2,4-Trichlorobenzene	.4	U	Pyrene	.4	Ū
Naphthalene	.4	U	Retene	.4	U
4-Chloroaniline	.4	Ū	Butylbenzylphthalate	.025	${f J}$
Hexachlorobutadiene	.4	U	Benzo(a)anthracene	.4	Ŭ
4-Chloro-3-Methylphenol	.4	U	3,3'-Dichlorobenzidine	.8	U
2-Methylnaphthalene	.4	Ū	Chrysene	.4	U
1-Methylnaphthalene	.4	Ū	Bis(2-Ethylhexyl) Phthalate	.044	f J
Hexachlorocyclopentadiene	.4	Ŭ	Di-N-Octyl Phthalate	.4	บ
2,4,6-Trichlorophenol	.4	Ŭ	Benzo(b)fluoranthene	.4	Ū
2,4,5-Trichlorophenol	.4	Ū	Benzo(k)fluoranthene	.4	Ŭ
2-Chloronaphthalene	.4	Ŭ	Benzo(a)pyrene	.4	Ū
2-Nitroaniline	.4	Ŭ	3B-Coprostanol	.8	Ŭ
Dimethylphthalate	.4	Ŭ	Indeno(1,2,3-cd)pyrene	.4	Ŭ
2,6-Dinitrotoluene	.4	Ŭ	Dibenzo(a,h)anthracene	.4	Ŭ
Acenaphthylene	.4	Ŭ	Benzo(ghi)perylene	.4	Ŭ
3-Nitroaniline	.4	Ŭ		• •	_
		_			

Authorized By: __

Release Date: 2/5/99

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Analysis Report for

Base/Neutral/Acids

Project Name: Hamilton Labree Rds. - 04 LIMS Project ID: 1096-99

Method: SW8270

Lab ID: OBW9027B2
QC Type: Laboratory Method Blank
Project Officer: Pam Marti

Date Prepared: 01/27/99 Matrix: Water Date Analyzed: 01/27/99

Units:

ug/L

Surrogate Recoveries

2-Fluorophenol	67	%
D5-Phenol	51	%
D4-2-Chlorophenol	82	%
1,2-Dichlorobenzene-D4	66	%
D5-Nitrobenzene	85	%
2-Fluorobiphenyl	78	%
Pyrene-D10	115	%
Terphenyl-D14	111	- %

Authorized By:

Release Date: 2/5/59

Page:

7411 Beach Dr E, Port Orchard Washington 98366

CASE NARRATIVE

February 4, 1998

Subject:

Hamilton/Labree Rds. Project

Sample(s):

99048005

Officer(s):

Pam Marti

By:

Bob Carrell

Organics Analysis Unit

ACID HERBICIDE ANALYSIS

ANALYTICAL METHOD(S): (Draft EPA Method 8085)

The water sample for acid herbicides was extracted following Manchester Laboratory's standard operating procedure for the extraction of herbicides. The herbicide samples were hydrolyzed at pH > 12, extracted with methylene chloride at pH < 2, solvent exchanged and derivatized along with two method blanks. These extracts were analyzed by capillary Gas Chromatography and Atomic Emission Detection (GC/AED). Confirmation of herbicides is performed by Gas Chromatography and Ion-Trap mass spectrometry (GC/ITD) or comparisons of elemental ratios of hetero-atoms to empirical formulas.

The method utilizes compound independent calibration (CIC) for quantitation of detected compounds. A calibration validation is performed each time CIC is used for target compounds. This is done by comparison of CIC to a single point calibration (SPC) of the target analyte being quantitated.

All analytes have a respective practical quantitation limit (PQL) that is higher than the corresponding method detection limit (MDL). If a target analyte is detected and its identification is unambiguously confirmed at a concentration below its PQL, the reported concentration is qualified as an estimate, 'J' qualifier.

BLANKS:

No target compounds were detected in the laboratory blanks, thus demonstrating that the system was free from contamination.

HOLDING TIMES:

All samples were extracted and analyzed within the method holding times.

SURROGATES:

The 2,4,6-tribromophenol and 2,4-dichlorophenylacetic acid surrogate recoveries were acceptable, ranging from 88% to 96% and 45% to 72% respectively.

MATRIX SPIKING:

Not applicable.

COMMENTS:

One non-target chlorinated compound, tentatively identified by GC/ITD as 2-propanol, 1,3-dichloro-, phosphate (3:1), was observed in the sample. This compound is used as a flame retardant (formerly used in children's sleepware) and was found to have an estimated quantity of 0.26 ug/L.

The target analytes picloram and dinoseb received the 'UJ' qualifier because we traditionally experience highly variable recoveries for these compounds.

The data is useable as qualified.

DATA QUALIFIER CODES

*		
U	-	The analyte was not detected at or above the reported result.
J	-	The analyte was positively identified. The associated numerical result is an <u>estimate</u> .
UJ	-	The analyte was not detected at or above the reported estimated result.
REJ	-	The data are <u>unusable</u> for all purposes.
NAF	-	Not analyzed for.
N	-	For organic analytes there is evidence the analyte is present in this sample.
NJ	-	There is evidence that the analyte is present. The associated numerical result is an estimate.
NC	-	Not Calculated
E	-	This qualifier is used when the concentration of the associated value exceeds the known calibration range.

Department of Ecology

Analysis Report for

Chlorophenoxy Herbicides

Project Name: Hamilton Labree Rds. - 04 LIMS Project ID: 1096-99

sample: 99048005

Date Collected: 01/26/99

Method: SW8085

Field ID: HL-5

Project Officer: Pam Marti

Date Prepared:	01/27/99	Matrix:	Water
Date Analyzed:	01/29/99	Units:	ug/L

\nalyte	Result	Qualifier	
A & Tricklerenhand	0.012	**	
2,4,6-Trichlorophenol	0.012	U U	
3,5-Dichlorobenzoic Acid	0.020		
+-Nitrophenol	0.035	U	
2,4,5-Trichlorophenol	0.012	U	
Dicamba I	0.020	U	
2.3,4,6-Tetrachlorophenol	0.011	U	
MCPP (Mecoprop)	0.040	ŭ	·
MCPA	0.040	ũ	<u>.</u>
Dichlorprop	0.022	Ũ	
3romoxynil	0.020	Ü	
2,4-D	0.020	ñ	
2,3,4,5-Tetrachlorophenol	0.011	ũ	
Trichlopyr	0.017	ũ	
Pentachlorophenol	0.010	U	
2,4,5-TP (Silvex)	0.016	U	
2,4,5-T	0.016	U	
2,4-DB	0.024	U	
Dinoseb	0.030	ហ	4 L
Bentazon	0.030	ū	
[oxynil Pinton	0.020	U	
Picloram	0.020	ŢĴ	-
Dacthal (DCPA)	0.016	U	
2,4,5-TB	0.018	ŭ	
Acifluorfen (Blazer)	0.079	ū	
Diclofop-Methyl	0.030	U	
Surrogate Recoveries			
2,4,6-Tribromophenol	96	%	The second control of
2,4-Dichlorophenylacetic acid	72	%	·

2,4,6-Tribromophenol	96	%
2,4-Dichlorophenylacetic acid	72	%

Authorized By: 3

Release Date: 2-4-69

Department of Ecology

Analysis Report for

Chlorophenoxy Herbicides

Project Name:

Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Lab ID:

OBW9027A1H

Method: SW8085

QC Type: Laboratory Method Blank Project Officer: Pam Marti

Date Prepared: 01/27/99

Matrix: Water

Date Analyzed: 01/29/99

Units: ug/L

Analyte	Result	Oualifier	
(ALERI) CO	Moult	Anamici	
2,4,6-Trichlorophenol	0.013	U	
3,5-Dichlorobenzoic Acid	0.021	Ŭ	
4-Nitrophenol	0.036	Ū	
2,4,5-Trichlorophenol	0.013	Ū	
Dicamba I	0.021	U	
2,3,4,6-Tetrachlorophenol	0.012	U	
MCPP (Mecoprop)	0.042	U	
MCPA	0.042	Ū	
Dichlorprop	0.023	Ü	
Bromoxynil	0.021	Ū	
2,4-D	0.021	Ū	
2,3,4,5-Tetrachlorophenol	0.012	Ū	
Trichlopyr	0.018	Ū	
Pentachlorophenol	0.010	Ū	
2,4,5-TP (Silvex)	0.017	Ū	
2,4,5-T	0.017	Ū	
2,4-DB	0.025	Ū	
Dinoseb	0.031	ŪJ	
Bentazon	0.031	Ü	
Ioxynil	0.021	Ū	
Picloram	0.021	ŪJ	
Dacthal (DCPA)	0.017	Ū	
2,4,5-TB	0.019	Ū	
Acifluorfen (Blazer)	0.083	Ŭ	
Diclofop-Methyl	0.031	Ŭ	
Surrogate Recoveries			
2,4,6-Tribromophenol	90	%	
2,4-Dichlorophenylacetic acid	45	%	

Authorized By: __

Release Date: 2-4-79

Department of Ecology

Analysis Report for

Chlorophenoxy Herbicides

Hamilton Labree Rds. - 04 roject Name:

LIMS Project ID: 1096-99

OBW9027A2H ab ID:

C Type: Laboratory Method Blank roject Officer: Pam Marti Date Prepared: 01/27/99 Method: SW8085 Matrix: Water

Date Analyzed: 01/29/99 Units:

ug/L

·									
\nalyte	Result	Qualifier	•		· · ·				
,4,6-Trichlorophenol	0.013	U							
,5-Dichlorobenzoic Acid	0.021	Ŭ							
-Nitrophenol	0.036	Ū				-			
.,4,5-Trichlorophenol	0.013	Ū							
Dicamba I	0.021	Ū				5			
1,3,4,6-Tetrachlorophenol	0.012	U							
ACPP (Mecoprop)	0.042	U							
MCPA	0.042	U							
Dichlorprop	0.023	U							
3romoxynil	0.021	U			-			÷=	
!,4-D	0.021	U .							
2,3,4,5-Tetrachlorophenol	0.012	Ü							
[richlopyr	0.018	\mathbf{U}			•				
Pentachlorophenol	0.010	U							
!,4,5-TP (Silvex)	0.017	U		•					
?,4,5-T	0.017	U							
2,4-DB	0.025	U							
Dinoseb	0.031	UJ							
3entazon	0.031	\mathbf{U}							
.oxynil	0.021	U							
Picloram	0.021	UJ							
Dacthal (DCPA)	0.017	U							
2,4,5-TB	0.019	U							
Acifluorfen (Blazer)	0.083	U					•		
Diclofop-Methyl	0.031	U							
Surrogate Recoveries									
2,4,6-Tribromophenol	88	%	\neg	•				-	
2,4-Dichlorophenylacetic acid	51	%							
				. w Y	5 min man 51			~	

uthorized By: 132

Release Date: 2-4-99

7411 Beach Dr E, Port Orchard Washington 98366

CASE NARRATIVE

February 10, 1999

Subject:

Hamilton Labree Rds Project: Water Sample

Samples:

99048005

Officer(s):

Pam Marti

By:

Norman Olson (CC)
Organics Analysis Unit

Hetero-atom Containing Compound Screen (AED SCREEN)

ANALYTICAL METHOD(S):

The water sample was extracted following Manchester Laboratory's standard operating procedure for the extraction of pesticides from water. The sample was extracted with methylene chloride followed by solvent exchange to iso-octane. The analytical method used was proposed EPA SW-846 Method 8085. The extract was analyzed using capillary gas chromatography and atomic emission detection (GC/AED) to screen the samples for hetero-atom containing compounds. A gas chromatography and ion trap mass spectrometry (GC/MS) system was used to obtain identification and/or confirmation for such detected compounds if warrented.

With this procedure, Method 8085 in screening mode, all organic compounds with the following characteristics would have been detected if present in the sample.

- 1. Contains one or more of the hetero-atoms: Cl, Br, I, N, S or P
- 2. Is semi-volatile, extractable and chromatographical
- 3. Is at sufficient concentration (low to high ppt range)

This analytical method used for the water sample was limited to a minimal level of quality assurance for quantitation. All compounds reported present have been unambiguously confirmed to be present, and the associated quantitation is qualified as estimated. Furthermore, all target analytes reported as not detected have practical quantitation limits (PQLs) that are qualified 'UJ' as estimates.

BLANKS:

None of the target compounds were detected in the method blanks. Hence, the method blanks demonstrate the system was free from contamination.

SURROGATE(S):

Recoveries were acceptable.

COMMENTS:

No hetero-atom containing compounds were detected in the sample.

The data is useable as qualified.

DATA QUALIFIER CODES

U	-	The analyte was not detected at or above the reported result.
J	-	The analyte was positively identified. The associated numerical result is an estimate.
UJ	-	The analyte was not detected at or above the reported estimated result.
N	-	For organic analytes there is evidence the analyte is present in this sample.
NJ	-	There is evidence that the analyte is present. The associated numerical result is an estimate.

Department of Ecology

Analysis Report for

Chlorinated Pesticides (GC/AED)

Project Name: Hamilton Labra	ee Rds	04	LIMS Project ID: 1096-99
Sample: 99048005 Field ID: HL-5 Project Officer: Pam Marti		Date Pr	ollected: 01/26/99 Method: SW8085 repared: 01/28/99 Matrix: Water nalyzed: 02/05/99 Units: ug/L
Analyte	Result	Qualifier	
Alpha-BHC	0.010	υJ	Surrogate Recoveries
Beta-BHC	0.010	IJ	
Gamma-BHC (Lindane)	0.010	UJ	4,4-Dibromooctafluorobiphenyl 74 %
Delta-BHC	0.010	IJ	Decachlorobiphenyl 101 %
Heptachlor	0.010	$\mathbf{U}\mathbf{J}$	
Aldrin	0.010	$\mathbf{U}\mathbf{I}$	•
Heptachlor Epoxide	0.010	$\mathbf{U}\mathbf{J}$	
Trans-Chlordane (Gamma)	0.010	$\mathbf{U}\mathbf{J}$	
Endosulfan I	0.010	UJ	,
Dieldrin	0.010	$\mathbf{U}\mathbf{J}$	••
4,4'-DDE	0.010	UJ	·
Endrin	0.010	UJ	
Endosulfan II	0.010	UJ	,
4,4'-DDD	0.010	UJ	
Endrin Aldehyde	0.010	\mathbf{u}	
Endosulfan Sulfate	0.010	$\overline{\mathbf{n}}$	
4,4'-DDT	0.010	$\mathbf{U}\mathbf{J}$	
Endrin Ketone	0.010	U J	•
Methoxychlor	0.010	$\overline{\Omega}$	
Alpha-Chlordene	0.010	UJ	
Gamma-Chlordene	0.010	$\mathbf{U}\mathbf{J}$	
Oxychlordane	0.010	UJ	· · · · · · · · · · · · · · · · · · ·
DDMU	0.010	UJ	- · · · · · · · · · · · · · · · · · · ·
Cis-Chlordane (Alpha-Chlordane	0.010	UJ	
Cis-Nonachlor	0.010	$\mathbf{u}_{\mathbf{l}}$	Tr.
Kelthane	0.040	$\mathbf{U}\mathbf{J}$	
Captan	0.027	$\mathbf{U}\mathbf{J}$	
2,4'-DDE	0.010	Πī	
Trans-Nonachlor	0.010	ŪJ	
2,4'-DDD	0.010	<u>UJ</u>	
2,4'-DDT	0.010	U J	
Captafol	0.050	ប្រ	
Mirex	0.010	UJ	
Toxaphene	0.30	UJ	
Hexachlorobenzene	0.010	UJ	
Pentachloroanisole	0.010	Ωĵ	
PCB - 1254	0.063	IJ	
PCB - 1260	0.063	UJ	
		<u>,</u>	

Release Date:

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Authorized By:

Department of Ecology

Analysis Report for

Chlorinated Pesticides (GC/AED)

Project Name: Hamilton Labree Rds. - 04 LIMS Project ID: 1096-99

Lab ID: OBW9028B1 Method: SW8085

QC Type: Laboratory Method Blank
Project Officer: Pam Marti

Date Prepared: 01/28/99 Matrix: Water
Date Analyzed: 02/05/99 Units: ug/L

Analyte	Result	Qualifier			
Alpha-BHC	0.010	UJ	Surrogate Recoveries		
Beta-BHC	0.010	$\mathbf{u}_{\mathbf{l}}$			
Gamma-BHC (Lindane)	0.010	UJ	4,4-Dibromooctafluorobiphenyl	84	%
Delta-BHC	0.010	UJ	Decachlorobiphenyl	106	%
Heptachlor	0.010	IJ			
Aldrin	0.010	UJ			
Heptachlor Epoxide	0.010	UJ			
Trans-Chlordane (Gamma)	0.010	UJ			
Endosulfan I	0.010	បា	-		
Dieldrin	0.010	UJ			
4,4'-DDE	0.010	UJ			
Endrin	0.010	UJ			
Endosulfan II	0.010	$\mathbf{U}\mathbf{J}$			
4,4'-DDD	0.010	UJ			
Endrin Aldehyde	0.010	UJ			
Endosulfan Sulfate	0.010	UJ	•		
4,4'-DDT	0.010	υJ			
Endrin Ketone	0.010	UJ			
Methoxychlor	0.010	ŨĴ			
Alpha-Chlordene	0.010	ni ni			
Gamma-Chlordene	0.010	ប្ប			
Oxychlordane	0.010	UJ			
DDMU	0.010	<u>ni</u>			
Cis-Chlordane (Alpha-Chlordane	0.010	Ū J			
Cis-Nonachlor	0.010	ŪJ			
Kelthane	0.040	UJ			
Captan	0.027	UJ			
2,4'-DDE	0.010	UJ			
Trans-Nonachlor	0.010	nı			
2,4'-DDD	0.010	ŪJ			
2,4'-DDT	0.010	UJ			
Captafol	0.050	UJ			
Mirex	0.010	<u>ni</u>			
Toxaphene	0.30	Ωĵ			
Hexachlorobenzene	0.010				
Pentachloroanisole	0.010				
PCB - 1254	0.063				
PCB - 1260	0.063	UJ			
		·			

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Department of Ecology

Analysis Report for

Organophosphorous Pesticides (GC/AED)

Project Name: Hamilton Labree Rds 04				LIMS Project ID: 1096-99			
Sample: 99048005 Field ID: HL-5 Project Officer: Pam Marti		Date Pr	ollected: 01/26/99 repared: 01/28/99 nalyzed: 02/05/99	Method: Matrix: Units:		-	
Analyte	Result	Qualifier					
Demeton-O	0.014	UJ	Surrogate Recove	ries			
Sulfotepp	0.012	ŪĴ				- · · -	
Demeton-S	0.014	UJ	Triphenyl Phosph	ate	112		%
Fonofos	0.012	UJ					
Disulfoton (Di-Syston)	0.012	UJ		•			
Methyl Chlorpyrifos	0.016	UJ					
Fenitrothion	0.014	ŬĴ	,				
Malathion	0.016	Ŭ J					
Chlorpyriphos	0.016	ŬĴ					-
Merphos (1 & 2)	0.024	ŬĴ					
Ethion	0.014	ŬĴ					
Carbophenothion	0.020	ŬĴ					
EPN	0.020	ប័រ					
Azinphos Ethyl	0.032	ŬĴ		-	-		
Ethoprop	0.016	ŬĴ					
Phorate	0.014	ŬĴ					
Dimethoate	0.016	ŭi					
Diazinon	0.016	ŬĴ					
Methyl Parathion	0.014	Ü					
Ronnel	0.014	Ü					
Fenthion	0.014	Ü			-		
Parathion	0.014	U J			-		
Fensulfothion	0.020	UJ					
Bolstar (Sulprofos)	0.020	Ü					
Imidan	0.014	UJ					
				-			
Azinphos (Guthion)	0.032	UJ					
Coumaphos	0.024	IJ					
Dichlorvos (DDVP)	0.016	UJ	i				
Mevinphos	0.020	UJ					
Dioxathion	0.034	UJ			-		
Propetamphos Mathyl Pamoyon	0.040	\boldsymbol{u}					
Methyl Paraoxon	0.036	Ų					
Phosphamidan Tetrophomicanhon (Cordons)	0.048	UJ					
Tetrachlorvinphos (Gardona)	0.040	UJ			-		
Fenamiphos Tributes (DEE)	0.030	UJ					
Tribufos (DEF)	0.028	UJ		-			
Abate (Temephos)	0.12	UJ					
							_ =====
					-		
Authorized By:			Release Date:		P	age:	1

Department of Ecology

Analysis Report for

Organophosphorous Pesticides (GC/AED)

Project Name: LIMS Project ID: 1096-99 Hamilton Labree Rds. - 04

Method: SW8085

Lab ID: OBW9028B1
QC Type: Laboratory Method Blank
Project Officer: Pam Marti Date Prepared: 01/28/99 Matrix: Water Date Analyzed: 02/05/99 Units: ug/L

Analyte	Result	Qualifier		_
Demeton-O	0.014	UJ	Surrogate Recoveries	
Sulfotepp	0.012	UJ		_
Demeton-S	0.014	$\mathbf{U}\mathbf{J}$	Triphenyl Phosphate 119 %]
Fonofos	0.012	UJ		-
Disulfoton (Di-Syston)	0.012	UJ		
Methyl Chlorpyrifos	0.016	UJ		
Fenitrothion	0.014	UJ		
Malathion	0.016	UJ		
Chlorpyriphos	0.016	UJ		
Merphos (1 & 2)	0.024	\boldsymbol{u}		
Ethion	0.014			
Carbophenothion	0.020	Πl		
EPN	0.020	UJ	•	
Azinphos Ethyl	0.032	UJ		
Ethoprop	0.016	UJ		
Phorate	0.014	$\mathbf{U}\mathbf{J}$		
Dimethoate	0.016	$\mathbf{U}\mathbf{J}$		
Diazinon	0.016	U J		
Methyl Parathion	0.014	UJ		
Ronnel	0.014	UJ		
Fenthion	0.014	UJ		
Parathion	0.016	U J		
Fensulfothion	0.020	UJ		
Bolstar (Sulprofos)	0.014	UJ		
Imidan	0.022	UJ		
Azinphos (Guthion)	0.032	UJ		
Coumaphos	0.024	UJ		
Dichlorvos (DDVP)	0.016	UJ		
Mevinphos	0.020	U1		
Dioxathion	0.034	UJ		
Propetamphos	0.040	UJ	·	
Methyl Paraoxon	0.036	UJ		
Phosphamidan	0.048	UJ		
Tetrachlorvinphos (Gardona)	0.040			
Fenamiphos	0.030			
Tribufos (DEF)	0.028			
Abate (Temephos)	0.12	UJ		
_				

Authorized By:	Release Date:	Page:	1
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Department of Ecology

Analysis Report for

Nitrogen Containing Pesticides

Project Name: Hamilton Labree Rds. - 04 LIMS Project ID: 1096-99

Sample: 99048005

Field ID: HL-5

Project Officer: Pam Marti

Date Collected: 01/26/99

Method: SW8085

Date Prepared: 01/28/99

Matrix: Water

Date Analyzed: 02/05/99

Units: ug/L

Analyte	Result	Qualifier	Analyte	Result	Qualifier
Dichlobenil	0.040	IJ	Fenarimol	0.060	tu
Tebuthiuron	0.030	ŬĴ	Diuron	0.12	Ū
Propachlor (Ramrod)	0.048	Ωĵ	Di-allate (Avadex)	0.14	ŬJ .
Ethalfluralin (Sonalan)	0.030	ŪJ ·	Profluralin	0.048	Ū
Treflan (Trifluralin)	0.030	IJ	Metalaxyl	0.12	UJ -
Simazine	0.020	Πl	Cyanazine	0.030	υj
Atrazine	0.020	UJ	- 3		
Pronamide (Kerb)	0.079	ŪJ	Surrogate Recoveries		1
Terbacil	0.060	ŪĴ			
Metribuzin	0.020	ŪJ	1,3-Dimethyl-2-nitrobenzene	120	%
Alachlor	0.071	ŢŢ			
Prometryn	0.020	UJ			
Bromacil	0.079	UJ			
Metolachlor	0.079	IJ			
Diphenamid	0.060	Πl			
Pendimethalin	0.030	UJ	_		
Napropamide	0.060	UJ			
Oxyfluorfen	0.079	. UJ			
Norflurazon	0.040	$\mathbf{U}\mathbf{J}$			
Eptam	0.040	$\mathbf{U}\mathbf{J}$			
Butylate	0.040	$\mathbf{U}\mathbf{J}$			
Vernolate	0.040	UJ			
Cycloate	0.040	UJ			
Benefin	0.030	UJ	•		
Prometon (Pramitol 5p)	0.020	UJ			•
Propazine `	0.020	UJ			
Chlorothalonil (Daconil)	0.048	IJ			
Triallate	0.060	UJ			
Ametryn	0.020	UJ			
Terbutryn (Igran)	0.020	IJ			
Hexazinone	0.030	Πl	_ # . #		
Pebulate	0.040	UJ	1		
Molinate	0.040	ŪJ	· -		
Chlorpropham	0.079	ŬĴ			
Atraton	0.030	ŪĴ			
Triadimefon	0.052	ŬĴ			
MGK264	0.16	Ü	•		
Butachlor	0.12	ŬĴ			

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Department of Ecology

Analysis Report for

Nitrogen Containing Pesticides

Project Name:

Authorized By: __

Hamilton

oree Rds. - 04

LIMS Project ID: 1096-99

Page:

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Method: SW8085 Matrix: Water

Lab ID: OBW9028B1
QC Type: Laboratory Method Blank
Project Officer: Pam Marti

Date Prepared: 01/28/99 Date Analyzed: 02/05/99

Units: ug/L

Dichlobenil 0.040	Analyte	Result	Qualifier	Analyte	Result	Qualifier
Propachior (Ramrod)	Dichlobenil		UJ	Fenarimol	0.060	
Ethalfluralin (Sonalan)	Tebuthiuron	0.030	UJ	Diuron	0.12	UJ I
Ethalfluralin (Sonalan)	Propachlor (Ramrod)			Di-allate (Avadex)	0.14	UI
Treflan (Trifluralin)	Ethalfluralin (Sonalan)	0.030	UJ	Profluralin	0.048	UJ
Atrazine 0.020 UJ Surrogate Recoveries Terbacil 0.060 UJ Metribuzin 0.020 UJ IAJachlor 0.071 UJ Prometryn 0.020 UJ Metolachlor 0.079 UJ Metolachlor 0.060 UJ Pendimethalin 0.030 UJ Pendimethalin 0.030 UJ Napropamide 0.060 UJ Oxyfluorfen 0.079 UJ Norflurazon 0.040 UJ Eptam 0.040 UJ Eptam 0.040 UJ Eptam 0.040 UJ Cycloate 0.040 UJ Vernolate 0.040 UJ Vernolate 0.040 UJ Propazine 0.020 UJ Trialiate 0.060 UJ Ametryn 0.020 UJ Terbutryn (Igran) 0.020 UJ Hexazinone 0.030 UJ Pebulate 0.079 UJ Atraton 0.030 UJ Prizidimefon 0.052 UJ Molinate 0.079 UJ Atraton 0.030 UJ Triadimefon 0.052 UJ Molinate 0.052 U	Treflan (Trifluralin)	0.030	UJ	Metalaxyl	0.12	UJ
Pronamide (Kerb)	Simazine	0.020	UJ	Cyanazine	0.030	UJ
Terbacil 0.060 UJ 1.3-Dimethyl-2-nitrobenzene 190 %				•		}
Metribuzin 0.020 UJ I,3-Dimethyl-2-uitrobenzene 190 % Alachlor 0.071 UJ UJ Prometryn 0.020 UJ Bromacil 0.079 UJ UJ Pendimethalin 0.079 UJ Diphenamid 0.060 UJ Pendimethalin 0.030 UJ Napropamide 0.060 UJ VI VI Norflurazon 0.040 UJ VI Spatam 0.040 UJ VI Vernolate 0.040 UJ VI Vernolate 0.040 UJ VI Vernolate 0.040 UJ VI Propazine 0.030 UJ VI Propazine 0.020 UJ VI Chlorothalonil (Daconil) 0.048 UJ Triallate 0.060 UJ Ametryn 0.020 UJ Hexazinone 0.030 UJ Pebulate 0.0.0 UJ	Pronamide (Kerb)			Surrogate Recoveries		
Alachlor Prometryn 0.020 UJ Bromacil 0.079 With Metolachlor 0.079 Uj Diphenamid 0.060 Uj Pendimethalin 0.030 Uj Napropamide 0.060 Uj Oxyfluorfen 0.079 Norflurazon Eptam 0.040 Uj Eptam 0.040 Uj Eytam 0.040 Uj Vernolate 0.040 Uj Vernolate 0.040 Uj Ebenefin 0.030 Uj Propazine 0.020 Uj Chlorothalonil (Daconil) Triallate 0.060 Uj Hexazinone 0.030 Uj Pebulate 0.030 Uj Pebulate 0.040 Uj Chlorpropham 0.020 Uj Chlorpropham 0.020 Uj Chlorpropham 0.030 Uj Chlorpropham 0.079 Uj Atraton 0.030 Uj Triadimefon 0.030 Uj Triadimefon 0.052 Uj MGK264 0.16					_	
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MGK264 0.16 UJ						
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§	Butacnior	0.12	O J			

Release Date:

Washington Department of Ecology

Manchester Environmental Laboratory 7411 Beach Drive East Port Orchard, WA 98366

February 10, 1999

TO:

Pam Marti

FROM:

Jim Ross, Manchester Lab

SUBJECT:

Quality Assurance memo for the Hamilton Labree (Wk 4) project

SUMMARY

Data for this project met all quality assurance and quality control criteria with no qualification necessary. The GFAA is not operating properly, so data for Pb, As, Se and Tl is not available at this time. These analytes were not detected at ICP levels (20-50 ug/L)

SAMPLE RECEIPT

The samples were received by the Manchester Laboratory on 01/27/99

HOLDING TIMES

All analyses were performed within the specified holding time.

INSTRUMENT CALIBRATION

Instrument calibration was performed before each analytical run and checked by initial calibration verification standards and blanks. The AA calibration curves returned a correlation coefficient of 0.995 or better. Continuing calibration standards and blanks were analyzed at a frequency of 10% during the run and again at the end of the analytical run. All initial and continuing calibration verification standards and blanks were within the relevant control limits.

PROCEDURAL BLANKS

The procedural blanks associated with these samples showed no significant level of analytes.

SPIKED SAMPLE ANALYSES

All spike and duplicate spike recoveries met the acceptance criteria (75-125%)

PRECISION DATA

Precision estimates based on duplicate spike analysis were all within the acceptance criteria for duplicate analysis $(\pm 20\%)$

LABORATORY CONTROL SAMPLE (LCS) ANALYSES

All LCS analyses were within the acceptance criteria for the individual analytes.

Please call Jim Ross at (360) 871-8808 to further discuss this project.

Department of Ecology

Analysis Report for

Inductively Coupled Plasma, Total Recoverable

Project Name:

Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Sample: 99048005 Date Collected: 01/26/99

Method: BPA200.7

Field ID: HL-5

Date Prepared: 02/01/99

Matrix: Water

Project Officer: Pam Marti

Date Analyzed: 02/08/99

Units: ug/L

Analyte	Result	Qualifier		
Antimony	30	U		
Arsenic	0.4	U		
Beryllium	1	U	=	
Cadmium	4	Ŭ		•
Chromium	5	Ü		
Copper	7.3	_		
Lead	1	U		-
Nickel	10	ŭ		
Selenium	1	Ŭ		
Silver	$\bar{3}$	Ŭ		
Thallium	0.2	Ŭ		
Zinc	5.7	•		

Authorized By:

Release Date: 2/12/97

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Department of Ecology

Analysis Report for

Inductively Coupled Plasma, Total Recoverable

Project Name:

Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Method: BPA200.7

Sample: 99048005 (Matrix Spike LMX2) Date Collected: 01/26/99 Field ID: HL-5 Date Prepared: 02/01/99

Matrix: Water

Project Officer: Pam Marti

Date Prepared: 02/01/99 Date Analyzed: 02/08/99

Units:

% Recovery

Analyte	Result	Qualifier
Antimony	102	
Arsenic	107	
Beryllium	94	
Cadmium	101	
Chromium	98	
Copper	98	
Lead	107	
Nickel	97	
Selenium	109	
Silver	103	
Thallium	102	
Zinc	98	
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Department of Ecology

Analysis Report for

Inductively Coupled Plasma, Total Recoverable

Project Name:

Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Sample: 99048005 (Matrix Spike LMX1) Date Collected: 01/26/99

Method: EPA200,7 Date Prepared: 02/01/99

Field ID: HL-5

Matrix: Water

Project Officer: Pam Marti

Date Analyzed: 02/08/99

Units: % Recovery

Result Qualifier Analyte 102 Antimony 110 Arsenic Beryllium 97 Cadmium 103 Chromium 100 Copper 101 Lead 108 Nickel 99 111 Selenium Silver 104 Thallium 103 Zinc 100

Authorized By:

Release Date: 2/12/55

Department of Ecology

Analysis Report for

Inductively Coupled Plasma, Total Recoverable

Project Name:

Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Lab ID: M9032WB1

Method: EPA200.7

Matrix: Water

QC Type: Laboratory Method Blank Project Officer: Pam Marti

Date Prepared: 02/01/99 Date Analyzed: 02/08/99

Units: ug/L

Analyte	Result	Qualifier
A -43	30	υ
Antimony		
Arsenic	0.4	U
Beryllium	1	${f U}$
Cadmium	4	U
Chromium	5	U
Copper	5	U
Lead	1	U
Nickel	10	U
Selenium	1	U
Silver	3	U
Thallium	0.2	U
Zinc	5	U

Authorized By:

Department of Ecology

Analysis Report for

Inductively Coupled Plasma, Total Recoverable

Project Name:

Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Method: EPA200.7

Lab ID: M9032WL2

Matrix: Water

QC Type: Laboratory Control Sample Project Officer: Pam Marti

Date Prepared: 02/01/99 Date Analyzed: 02/08/99

Units: ug/L

Analyte	Result	Qualifier
Antimony	101	%
Arsenic	106	%
Beryllium	90	%
Cadmium	99	%
Chromium	97	%
Copper	96	%
Lead	108	%
Nickel	97	%
Selenium	113	%
Silver	103	%
Thallium	103	%
Zinc	96	%

Authorized By:

Release Date: 2//2/99

Department of Ecology

Analysis Report for

Mercury

Project Name: Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Project Officer: Pam Marti **Date Reported:** 03-FEB-99

Method: EPA245.1 Matrix: Water

Analyte: Mercury

Sample	QC	Field ID	Result	Qualifier	Units	Collected	Analyzed
99048005 99048005 99048005 M9033WG M9033WH	Matrix		0.03 104 % 99 % 110 % 0.03	u u	ug/L	01/26/99 01/26/99 01/26/99	02/03/99 02/03/99 02/03/99 02/03/99 02/03/99

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Department of Ecology

Analysis Report for

Alkalinity as Carbonate

Project Name: Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Project Officer: Pam Marti Pate Reported: 29-JAN-99 Method: EPA310.1 Matrix: Water

Matrix:

Analyte: Alkalinity as Carbonate

Sample	QC Field ID	Result Qualific	er Units	Collected	Analyzed
99048005 99048005 99048105	HE-5 Duplicate Matrix Spiles	71.7 71.4 101 %	mg/L mg/L	01/26/99 01/26/99 01/26/99	01/29/98 01/29/98 01/29/98

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Department of Ecology

Analysis Report for

Alkalinity as Bicarbonate

Project Name Pamilton Labree Rds. - 04

LIMS Project ID: 1096-99

! roject Offic r: him Marti Date Reporte i: 0-JAN-09 Method: EPA310.2

Matrix: Water

Analyte: Alkalinity as Bicarbonate

Sample QC Ted ID	Result Qualifier	Units	Collected	Analyzed
99048(05 FL-5	0.0	mg/L	01/26/99	01/29/98
99048(05 Duplica:		mg/L	01/26/99	01/29/98

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Department of Ecology

Analysis Report for

Hardness

Project Name: Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Project Officer: Pam Marti Date Poported: 10-FEB-99

Method: SM2340B Matrix: Water

Analyte: Hardness

Sample	QC	Field ID	Result	Qualifier	Units	Collected	Analyzed
99048005 99048005 M9039HB M9039HL		HL-5	118 118 0.2 95 %	U	mg/L .mg/L mg/L	01/26/99 01/26/99	02/08/98 02/08/98 02/08/98 02/08/98

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Department of Ecology

Analysis Report for

Total Dissolved Solids

Project Name: Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Project Officer: Pam Marti

Method:

EPA160.1 Water

Date Reported: 01-FEB-99

Matrix:

Total Dissolved Solids Analyte:

Sample	QC Field ID	Result	Qualifier Units	Collected	Analyzed
99048005	HL-5	212	mg/L	01/26/99	01/28/99
	Duplicate	210	mg/L	01/26/99	01/28/99

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Department of Ecology

Analysis Report for

Nitrite/Nitrate

Project Name: Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Project Officer: Pam Marti **Date Reported:** 01-FEB-99

Method: EPA353.2

Matrix: Water

Analyte: Nitrite-Nitrate

Sample	QC Field ID	Result Qualifier	Units	Collected	Analyzed
99048005 99048005 99048005	HL-5 Duplicate Matrix Spike	2.50 2.51 90.2 %	mg/L mg/L	01/26/99 01/26/99 01/26/99	01/27/99

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Department of Ecology

Analysis Report for

Chlorides

Project Name: Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Project Officer: Pam Marti Date Reported: 29-JAN-99 Method: EPA300.0 Matrix: Water Analyte: Chloride

Sample	QC_	Field ID	Result	Qualifier U	nits	Collected	Analyzed
99048005 99048005 99048005			44.4 43.8 95.0 %	m	g/L g/L	01/26/99 01/26/99 01/26/99	01/27/99 01/27/99 01/27/99

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Department of Ecology

Analysis Report for

Sulfate

Project Name: Hamilton Labree Rds. - 04

LIMS Project ID: 1096-99

Project Officer: Pam Marti Date Reported: 29-JAN-99 Method: EPA300.0 Matrix: Water

Analyte: Sulfate

Sample	QC Field ID	Result Qua	lifier Units	Collected	Analyzed
99048005 99048005 99048005	HL-5 Duplicate Matrix Spike	0.885 0.905 86.1 %	mg/L mg/L	01/26/99 01/26/99 01/26/99	01/27/99 01/27/99 01/27/99

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